

Data Path : Z:\HPCHEM1\BNA N\DATA\BN031318\
 Data File : BN000363.D
 Acq On : 13 Mar 2018 23:19
 Operator : JU/SJ
 Sample : SSTDCCC020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD02001

Quant Time: Mar 14 08:15:40 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN030918MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 14 08:11:44 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	197855	20.00	ng/ul	0.00
2) Naphthalene-d8	11.29	136	936363	20.00	ng/ul	0.00
6) Acenaphthene-d10	15.06	164	537586	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.79	188	1227142	20.00	ng/ul	0.00
17) Chrysene-d12	21.90	240	1231152	20.00	ng/ul	0.00
22) Perylene-d12	24.39	264	1141238	20.00	ng/ul	0.00
System Monitoring Compounds						
7) Acenaphthylene-d8	14.76	160	1129162	21.16	ng/ul	0.00
10) Fluorene-d10	16.05	176	751063	20.79	ng/ul	0.00
14) Anthracene-d10	17.89	188	1201009	20.95	ng/ul	0.00
18) Pyrene-d10	20.14	212	1347495	22.37	ng/ul	0.00
25) Benzo(a)pyrene-d12	24.23	264	1087536	20.87	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.34	128	999179	20.221	ng/ul	99
4) 2-Methylnaphthalene	12.92	142	691512	19.881	ng/ul	98
5) 1-Methylnaphthalene	13.13	142	694489	19.896	ng/ul#	93
8) Acenaphthylene	14.79	152	1119492	20.879	ng/ul	98
9) Acenaphthene	15.13	153	759979	20.365	ng/ul	95
11) Fluorene	16.10	166	859485	20.882	ng/ul	93
13) Phenanthrene	17.83	178	1429858	20.317	ng/ul	100
15) Anthracene	17.92	178	1482947	20.767	ng/ul	99
16) Fluoranthene	19.80	202	1708372	22.467	ng/ul#	92
19) Pyrene	20.17	202	1724157	22.182	ng/ul#	91
20) Benzo(a)anthracene	21.89	228	1614760	21.098	ng/ul	98
21) Chrysene	21.95	228	1471693	20.262	ng/ul	100
23) Benzo(b)fluoranthene	23.63	252	1452260	21.340	ng/ul#	96
24) Benzo(k)fluoranthene	23.68	252	1424964	22.370	ng/ul#	96
26) Benzo(a)pyrene	24.28	252	1343910	20.750	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	26.96	276	1489709	21.332	ng/ul#	81
28) Dibenzo(a,h)anthracene	26.97	278	1242952	21.714	ng/ul#	93
29) Benzo(a,h,i)perylene	27.76	276	1231933	21.127	ng/ul#	89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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